

# Sebacic acid, but-3-enyl heptyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C21H38O4/c1-3-5-7-12-15-19-25-21(23)17-14-11-9-8-10-13-16-20(22)24-18-6

DNGGQHJUOLGNST-UHFFFAOYSA-N

C21H38O4

C=CCCOC(=O)CCCCCCCCC(=O)OCCCCCCC

354.52

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -254.06 | kJ/mol  | Joback Method  |
| hf            | -840.94 | kJ/mol  | Joback Method  |
| hfus          | 54.44   | kJ/mol  | Joback Method  |
| hvap          | 79.98   | kJ/mol  | Joback Method  |
| log10ws       | -6.19   |         | Crippen Method |
| logp          | 5.740   |         | Crippen Method |
| mcvol         | 317.330 | ml/mol  | McGowan Method |
| pc            | 1040.58 | kPa     | Joback Method  |
| rinpol        | 2472.00 |         | NIST Webbook   |
| rinpol        | 2472.00 |         | NIST Webbook   |
| tb            | 829.14  | K       | Joback Method  |
| tc            | 1016.39 | K       | Joback Method  |
| tf            | 468.99  | K       | Joback Method  |
| vc            | 1.240   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1002.79   | J/mol×K | 829.14          | Joback Method |
| cpg           | 1020.92   | J/mol×K | 860.35          | Joback Method |
| cpg           | 1037.99   | J/mol×K | 891.56          | Joback Method |
| cpg           | 1054.00   | J/mol×K | 922.77          | Joback Method |
| cpg           | 1068.99   | J/mol×K | 953.97          | Joback Method |
| cpg           | 1082.97   | J/mol×K | 985.18          | Joback Method |
| cpg           | 1095.98   | J/mol×K | 1016.39         | Joback Method |
| dvisc         | 0.0007636 | Pa×s    | 468.99          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003732 | Pa×s | 529.01 | Joback Method |
| dvisc | 0.0002110 | Pa×s | 589.04 | Joback Method |
| dvisc | 0.0001326 | Pa×s | 649.07 | Joback Method |
| dvisc | 0.0000902 | Pa×s | 709.09 | Joback Method |
| dvisc | 0.0000651 | Pa×s | 769.12 | Joback Method |
| dvisc | 0.0000493 | Pa×s | 829.14 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U356087&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U356087&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>rinpol:</b>  | Non-polar retention indices                     |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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